## Comment on "Bethe Ansatz Results for the 4f-Electron Spectra of a Degenerate Anderson Model"

In a recent letter [1], the author calculates the density of states for 4f electrons coupled to a conduction band in the framework of the Bethe ansatz (BA) solution for the degenerate Anderson model. It is claimed that the results qualitatively disagree with the results obtained for the same model but using a variational approach [2]. Even the high energy feature in the f-spectral function near the 4f-level energy  $\epsilon_f$ , i.e. the "normal" ionization peak (NIP), is argued to be qualitatively different in the two approaches. In the following we point out that this is not the case.

We concentrate on the  $U \to \infty$ , N-fold degenerate Anderson model. As the Bethe ansatz for this model can yield exact results only in the large bandwidth limit  $B \gg |\epsilon_f|$ , we confine our discussion to this limit. Zvyagin first presents BA results which show that the NIP is shifted towards the chemical potential from  $\epsilon_f$  in the Kondo limit  $n_f \approx 1$ , where  $n_f$  is the mean f-occupancy. He then states that the peak shifts in the opposite direction in our variational calculation [2]. That this is not the case is obvious from our results in the Kondo limit presented in Appendix C of reference [2]. This is also shown very clearly in Fig. 5 of our handbook article [3]. The NIP position is called  $\tilde{\epsilon}_f$  in Appendix C of reference [2]. For large N and the Kondo limit this real quantity is determined by the equation

$$\tilde{\epsilon}_f = \epsilon_f + Re\tilde{\Gamma}(\tilde{\epsilon}_f), \tag{1}$$

where

$$\tilde{\Gamma}(z) = N \int_{-B}^{0} d\epsilon |V(\epsilon)|^{2} / (z - \epsilon). \tag{2}$$

 $Re\tilde{\Gamma}(\epsilon)$  is positive for negative  $\epsilon$  with  $|\epsilon| \ll B$  (for a constant  $V(\epsilon)$  it is given by  $N\Delta \ln(B/|\epsilon|)$ ). It therefore follows without actually solving the equation that in our variational calculation the NIP shifts towards the chemical potential as in the BA solution. Zvyagin claims that the solution to this equation was found by us "in the complex form" [1]. Apparently he was confused by an unfortunate misprint in the second equality of our equation which determines  $\tilde{\epsilon}_f$ . It should read  $|\tilde{\epsilon}_f|$  instead of  $\tilde{\epsilon}_f$  in the argument of the logarithm. From the first part of the equation presented above it should be obvious that this is in fact a misprint.

We should mention that the results for the f-spectral function using our variational method [2] essentially agree with the low temperature solution of the NCA equations and numerical renormalization group results [4]. As noted by Zvyagin [1] other low energy characteristics obtained within our scheme qualitatively coincide with the BA ones.

In conclusion we have pointed out that in contrast to the claims in reference [1] our variational results for the high energy feature in the f-spectral function for the degenerate Anderson model, which has been used successfully in the description of photoemission spectra of Cerium compounds, qualitatively agree with the BA results.

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- [4] For a general discussion see: A. C. Hewson, *The Kondo problem to Heavy Fermions* (Cambridge University Press, Cambridge, 1993)